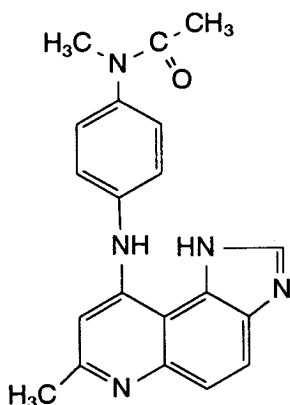


ACODAZOLE HYDROCHLORIDE

NSC - 305884



* HCl

Chemical Name:

N-Methyl-*N*-[4-[(7-methyl-1*H*-imidazo(4,5-*f*)quinolin-9-yl)amino]phenyl]acetamide, monohydrochloride

CAS Registry Number: 55435-65-9

Molecular Formula: C₂₀H₁₉N₅O · HCl

M.W. : 381.9

Approximate Solubility:

(mg/mL)

H₂O

> 50

pH 4 Acetate buffer

> 50

| | |
|-----------------------|---------|
| pH 9 Carbonate buffer | > 50 |
| 10% Ethanol | > 50 |
| Methanol | 10 - 12 |
| 95% Ethanol | 10 - 15 |
| 0.1 N HCl | > 50 |
| 0.1 N NaOH | < 1 |

Stability:

Bulk:

The bulk compound did not undergo significant decomposition when heated at 60 °C for 30 days in the dark (HPLC). The compound is quite hygroscopic.

Solution:

A solution (0.1 mg/mL) prepared in unbuffered water and stored at room temperature and exposed to light was stable for > 48 hours (HPLC).

Ultraviolet Absorption:

(Methanol)

| λ_{\max} | ϵ |
|------------------|-----------------|
| 345 \pm 2 nm | 15,300 - 16,600 |
| 316 \pm 2 nm | 12,400 - 13,200 |
| 255 \pm 2 nm | 27,900 - 30,500 |
| 227 \pm 2 nm | 21,100 - 22,400 |

High Performance Liquid Chromatography:

| | |
|---------------------------|--|
| Column: | μ -Bondapak C ₁₈ , 300 x 3.9 mm ID (Waters Assoc.) |
| Mobile Phase: | Methanol/water (containing 2% acetic acid), 40/60 (v/v) |
| Flow Rate: | 1.0 mL/min |
| Detection: | UV at 254 nm |
| Internal Standard: | Acetophenone, 1.28 μ g/mL |
| Retention Volume: | 25.0 mL (NSC-305884) 13.0 mL (I.S.) |

Toxicity Data:

Mouse(ip): LD₅₀: 161 mg/kg
NCI Screening Program Data Summary

Mouse(iv): LD₅₀: 87,590 μ g/kg
NCI Screening Program Data Summary